

## WEST Search History

DATE: Friday, July 18, 2003

**Set Name Query**

side by side

**Hit Count Set Name**

result set

*DB=USPT,PGPB; PLUR=YES; OP=OR*

L14	L13 or l12 or l6	68	L14
L13	L12 and mooren	4	L13
L12	L11 or l8	61	L12
L11	L10 and ocular	61	L11
L10	L9 and l1	206	L10
L9	L2 or l3	1779	L9
L8	L7 and l2	35	L8
L7	L5 and l1	58	L7
L6	L5 and mooren	11	L6
L5	L4 and inhibitor	357	L5
L4	L3 and ocular	361	L4
L3	cox near3 2	1533	L3
L2	celecoxib or deracoxib or valdecoxib or rofecoxib or etoricoxib	648	L2
L1	((514/403  514/406  514/473  514/316  514/474  514/378  514/318 )!..CCLS. )	4165	L1

END OF SEARCH HISTORY

=> e deracoxib/cn

E1 1 DERACOAT U 3051/CN  
E2 1 DERACOAT X 600W/CN  
E3 1 --> DERACOXIB/CN  
E4 1 DERAKANE/CN  
E5 1 DERAKANE 114/CN  
E6 1 DERAKANE 118/CN  
E7 1 DERAKANE 200/CN  
E8 1 DERAKANE 200, POLYMER WITH AMINOBENZOIC ACID HYDRAZIDE, 1,1'-  
-(METHYLENEDI-4,1-PHENYLENE) BIS (1H-PYRROLE-2,5-DIONE) AND 1,  
1'-(1-METHYLETHYLIDENE) BIS (4-(2-PROPENYLOXY) BENZENE)/CN  
E9 1 DERAKANE 200, POLYMER WITH N-ACETYL-N-(4-((4-(2,5-DIHYDRO-2,  
5-DIOXO-1H-PYRROL-1-YL) PHENYL) METHYL) PHENYL) ACETAMIDE AND 1,  
1'-(4-METHYL-1,3-PHENYLENE) BIS (1H-PYRROLE-2,5-DIONE)/CN  
E10 1 DERAKANE 200, POLYMER WITH N-ACETYL-N-(4-((4-(2,5-DIHYDRO-2,  
5-DIOXO-1H-PYRROL-1-YL) PHENYL) METHYL) PHENYL) ACETAMIDE, DOBEC  
KAN FT 3018, 1,1'-(4-METHYL-1,3-PHENYLENE) BIS (1H-PYRROLE-2,5  
-DIONE) AND 2,4,6-TR/CN  
E11 1 DERAKANE 350/CN  
E12 1 DERAKANE 411/CN

=> s e3

L1 1 DERACOXIB/CN

=> d l1

L1 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS

RN 169590-41-4 REGISTRY

CN Benzenesulfonamide, 4-[3-(difluoromethyl)-5-(3-fluoro-4-methoxyphenyl)-1H-  
pyrazol-1-yl]- (9CI) (CA INDEX NAME)

OTHER NAMES:

CN 4-[5-(3-Fluoro-4-methoxyphenyl)-3-(difluoromethyl)-1H-pyrazol-1-  
yl]benzenesulfonamide

CN **Deracoxib**

CN SC 046

CN SC 46

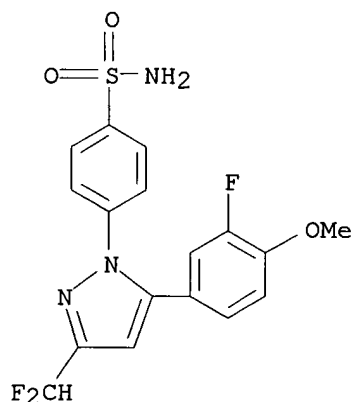
CN SC 59046

FS 3D CONCORD

MF C17 H14 F3 N3 O3 S

SR US Adopted Names Council

LC STN Files: ADISINSIGHT, CA, CAPLUS, TOXCENTER, USAN, USPATFULL, VETU



=> 514/403,406

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

41 REFERENCES IN FILE CA (1967 TO DATE)  
1 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA  
41 REFERENCES IN FILE CAPLUS (1967 TO DATE)

=> e rofecoxib/cn

E1	1	ROFANOL P 80/55/CN
E2	1	ROFANOL P 80/85/CN
E3	1 -->	ROFECOXIB/CN
E4	1	ROFELODINE/CN
E5	1	ROFEN 240/CN
E6	1	ROFENAID/CN
E7	1	ROFENAID 40/CN
E8	1	ROFENON/CN
E9	1	ROFERON A/CN
E10	1	ROFEROSE ST/CN
E11	1	ROFETAN F/CN
E12	1	ROFETAN GOT/CN

=> s e3

L2 1 ROFECOXIB/CN

=> d l2

L2 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS  
RN 162011-90-7 REGISTRY  
CN 2(5H)-Furanone, 4-[4-(methylsulfonyl)phenyl]-3-phenyl- (9CI) (CA INDEX NAME)

OTHER NAMES:

CN 3-Phenyl-4-[4-(Methylsulfonyl)phenyl]-2(5H)-furanone

CN MK 0966

CN MK 966

CN **Rofecoxib**

CN Vioxx

FS 3D CONCORD

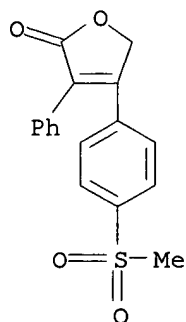
DR 186912-82-3

MF C17 H14 O4 S

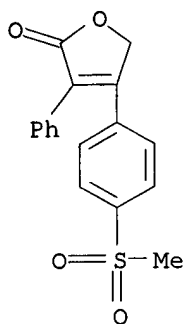
CI COM

SR CA

LC STN Files: ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, BIOSIS, BIOTECHNO, CA, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS, CIN, CSCHM, DIOGENES, DRUGNL, DRUGPAT, DRUGUPDATES, EMBASE, IPA, MRCK\*, PHAR, PHARMASEARCH, PROMT, RTECS\*, SYNTHLINE, TOXCENTER, USAN, USPAT2, USPATFULL  
(\*File contains numerically searchable property data)



⇒ 574/473, 474



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

299 REFERENCES IN FILE CA (1967 TO DATE)  
 9 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA  
 302 REFERENCES IN FILE CAPLUS (1967 TO DATE)

=> e valdecixib/cn

E1	1	VALCOAT 531/CN
E2	1	VALCONAZOLE/CN
E3	1 -->	VALDECOXIB/CN
E4	1	VALDET 4016/CN
E5	1	VALDET 79N/CN
E6	1	VALDETAMIDE/CN
E7	1	VALDIATE/CN
E8	1	VALDIBERINE/CN
E9	1	VALDIBERINE DIACETATE/CN
E10	1	VALDICE/CN
E11	1	VALDIPROMIDE/CN
E12	1	VALDISPERT/CN

=> s e3

L3 1 VALDECOXIB/CN

=> d 13

L3 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS

RN 181695-72-7 REGISTRY

CN Benzenesulfonamide, 4-(5-methyl-3-phenyl-4-isoxazolyl)- (9CI) (CA INDEX NAME)

OTHER NAMES:

CN 4-(5-Methyl-3-phenylisoxazol-4-yl)benzenesulfonamide

CN SC 65872

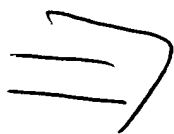
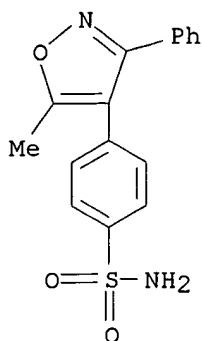
CN **Valdecixib**

FS 3D CONCORD

MF C16 H14 N2 O3 S

SR CA

LC STN Files: ADISINSIGHT, ANABSTR, BIOSIS, BIOTECHNO, CA, CAPLUS, CBNB, DRUGNL, DRUGPAT, DRUGUPDATES, EMBASE, IPA, PHAR, SYNTHLINE, TOXCENTER, USAN, USPAT2, USPATFULL



514/378

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

79 REFERENCES IN FILE CA (1967 TO DATE)

7 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

80 REFERENCES IN FILE CAPLUS (1967 TO DATE)

=> e etoricoxib/cn

E1	1	ETOPRINDOLE/CN
E2	1	ETOPRINE/CN
E3	1 -->	ETORICOXIB/CN
E4	1	ETORPHINE/CN
E5	1	ETORPHINE 3-HEMISUCCINATE/CN
E6	1	ETORPHINE DIHYDRATE/CN
E7	1	ETORPHINE HYDROCHLORIDE/CN
E8	1	ETORPHINE N-OXIDE/CN
E9	1	ETORPHINE-15,16-T2 HYDROCHLORIDE/CN
E10	1	ETORPHINE-15,16-T3/CN
E11	1	ETORPHINE-15-T HYDROCHLORIDE/CN
E12	1	ETORPHINE-3H/CN

=> s e3

L4 1 ETORICOXIB/CN

=> d 14

L4 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS

RN 202409-33-4 REGISTRY

CN 2,3'-Bipyridine, 5-chloro-6'-methyl-3-[4-(methanesulfonyl)phenyl]- (9CI)  
(CA INDEX NAME)

OTHER NAMES:

CN 5-Chloro-6'-methyl-3-[4-(methanesulfonyl)phenyl]-2,3'-bipyridine

CN **Etoricoxib**

CN MK 0663

CN MK 663

FS 3D CONCORD

MF C18 H15 Cl N2 O2 S

CI COM

SR CA

LC STN Files: ADISINSIGHT, BIOSIS, BIOTECHNO, CA, CAPLUS, CASREACT, DRUGNL,  
DRUGPAT, DRUGUPDATES, EMBASE, IPA, MEDLINE, PHAR, SYNTHLINE, TOXCENTER,  
USAN, USPAT2, USPATFULL